

CROSENTAL PROJECT CONTROL OF CONT 00213 5/126/60/009/04/004/033 18.1200 E032/E435 18. 2100 Garif'yanov, N.S. and Il'yasov. AUTHORS: Magnetic Resonance in Sodium Alloys TITLE: PERIODICAL: Fizika metallov i metallovedeniye, 1960, Vol 9, Nr pp 503-506 (USSR) Measurements are reported of the electron paramagnetic ABSTRACT: resonance in sodium and cesium alloys. The measurements were carried out on 300 and 9430 Mc/s at 295 and 90°K. The specimens were prepared in an argon atmosphere. In order to prevent the distortion of the lines due to the incomplete penetration of the high frequency field, the alloy was dispersed in paraffin, the particles being 4 µ in diameter (on the average). The sodium was 99.95% pure and contained about 0.04% of potassium. The width AH in the original sodium, measured as full width at half height, was 1.6 oe at 295°K and 9 oe at 90°K (Fig 1). These results are in agreement with those reported by Feher and Kip (Ref 3) and Gutowsky and Frank (Ref 4). Measurements on 300 Mc/s showed that AH for the electron paramagnetic is independent of the resonance curve for Na-Cs Card 1/5

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Magnetic Resonance in Sodium Alloys

concentration of cesium up to 0.5 at %. This is indicated in Fig 2. This result is in good agreement with Elliott's theory (Ref 5). Fig 2 shows the full width at half height of the electron paramagnetic width at half height of the electron paramagnetic resonance curve for different concentrations of sodium resonance curve for different concentrations of sodium and cesium. Measurements have also been made of the nuclear magnetic resonance in sodium alloys. The alloys investigated were Na-K., Na-Hg, and Na-Cs. The ratio of the Knight shift AH to the resonance value of the magnetic field H were measured. The results obtained are as follows:

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Magnetic Resonance in Sodium Alloys

Card 3/5

The Knight shift in sodium alloys (without correction for the chemical shift)

Alloy	Concentration of atoms in the alloy, %			△ H/H. ×	
	Na	<u>K</u>			
	100	0	3	0.1161	
Na-K	99.4	0.6		0.1160	
	93.8	6.2	- (0.1213	
	63	37 64	1	0.1287	
	36	64	- 1	0.1450	
	Na	Hg	1		
	97.2	2.8	1	0.1142	
Na-Hg	89.7	10.3		0.1122	
	85.9	14.1	1	0.1131	
	83.0	17.0		0.1149	
	81.0	19.0	i -1	0.1121	
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Magnetic Resonance	in Sodium	Alloys			
-	79	21	- 1	0.1152 0.1138	4
Na-Hg	77	23		0.1157	1
	75 70	25 30		0.1151	
	Na	Cs			
	99.7	0.3	1.	0.1125	
Na-Cs	99.5 99.4	0.5		0.1124	
na-cs	99.1	0.9		0.1152	
	98.7 94.2	1.3 5.8	· [0.1129	•
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E032/2435

Magnetic Resonance in Sodium Alloys

There are 3 figures, 1 table and 11 references,
4 of which are Soviet and 7 English.

ASSOCIATION: Fiziko-tekhnicheskiy institut Kazanskogo filiala AN SSSR (Physico-Technical Institute of the Kazan Branch AS USSR)

SUBMITTED: July 13, 1959

Card 5/5

IL'YASOV, A.V.; GARIF'YANOV, N.S.; RYZFMANOV, Yu.M.

Paramagnetic electron resonance in some types of natural cruds and in its heavy fractions. Ehim.i tekh.topl.i masel 6 no.1;28-31 Ja '61.

1. Figiko-tekhnicheskiy institut Kasanskogo filiala AW SSSR 1 Institut organichesköy khimii AW SSSR.

(Petroleum—Spectra)

23721:

8/057/61/031/006/008/019 B116/B203

24,7900(1144,1147,1163)

Garif'yanov, N. S., Il'yasov, A. V., and Ryzhmanov, Yu. M.

TITLE:

AUTHORS:

Electron paramagnetic resonance in some types of soot

PERIODICAL:

Zhurnal tekhnicheskoy fiziki, v. 31, no. 6, 1961, 694-698

TEXT: The authors studied the electron paramagnetic resonance (EPR) in various heat-treated types of soot at frequencies of V4 = 500 and V2 = 9450 Mc/sec in the temperature range of -193 to 300°C. They determined the relaxation times for gas conduit soot by the saturation method at 300 Mc/sec according to FOCT 785-49 TP 598 (GOST 785-49 column 598) as dependent on the temperature of the sample and the heat treatment. The measuring method used had been described earlier by N. S. Garif'yanov and B. M. Kozyrev (Ref. 3: ZhETF, 30, 272, 1956). The heat treatment consisted in heating to a certain temperature (maximum temperature 1200°C) without air access, with a holding time of 1 hr. The heat-treated soot was wetted with vaseline oil to eliminate the distortions on the EPR line at y and to obtain equal saturations of these lines on the whole sample

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Electron paramagnetic resonance in ... B116/B203

at V4. The samples were heated to 140°C (after heat treatment) without additional air removal by suction to eliminate the admorbed molecular oxygen. The authors investigated by the EPR method: furnace soot, gas conduit soot, and nozzle soot. Measurements were made in "oil" samples with sucked-off air at ν_1 = 300 and ν_2 = 9450 Mc/sec, and at -193, 20, and 300°C. A measurable effect was only observed with gas conduit soot. Results are tabulated. The authors found a strong dependence of the resonance line width AH on the temperature of heat treatment, and a weaker dependence on the temperature of the sample and on V. For all gas conduit soot samples, the splitting factor g was 2.003. The EPR curves show a Lorentz shape. The higher the heat treatment temperature and the temperature of the sample, the less the lines show saturation. The spinlattice relaxation time T, and the spin-spin relaxation time T, were determined at ν_1 by means of the saturation method for gas conduit soot samples from which the oxygen had been removed (Table 2). The strong concentration of paramagnetic centers, the equality of relaxation times $(T_1 \approx T_2)$, and the small line width in samples of gas conduit soot (in Card 2/6

8/057/61/031/006/008/019 B116/B203

Electron paramagnetic resonance in ...

heat treatment up to 900°C) suggest an exchange interaction between unpaired electrons. Evidently, the exchange is maintained also with a change in the temperature of the sample from -193 to 300°C, since also here $T_1 \approx T_2$. The Lorentz shape of the EPR curves also suggests an exchange interaction between paramagnetic centers. R. L. Collins, M. D. Bell and G. Kraus (Ref. 1: J. appl. phys., 30, 56, 1959) attempted to explain the rapid change of Δ H with increasing heat-treatment temperature up to 900-1000°C by the strong anisotropy of the g-factor. For anisotropic lines, the width of Δ H must depend very strongly on the frequency. The Δ H measured (Table 1)(at ν_1 and ν_2 differing by a factor of 30) differ only slightly. The data obtained confirm the assumption by J. Uebersfeld (Ref. 2: Ann. Phys., 13, 391, 1956). They explain the widening of the line by the reduction of T_1 due to the collision of unpaired electron with

carriers. The fact that no EPR were found with furnace soot and nozzle soot is explained by the circumstance that these types of soot are subjected to heating up to about 1200°C already during their formation. The temperature dependences of the lines in the gas conduit soot samples

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Electron paramagnetic resonance in ...

have not been explained so far. The presence of the adsorbed oxygen in ordinary and in heat-treated gas conduit soot samples reduces the relaxation times T₁ and T₂. The air is sucked off with difficulty from gas conduit soot samples exposed to air for a long time; therefore, the EPR line is wider in such samples as compared with fresh samples. There are 1 figure, 2 tables, and 9 references: 5 Soviet-bloc and 4 non-Soviet-bloc. The references to the English-language publications read as follows: N. Bloembergen, S. Wang. Phys. Rev., 23, 72, 1954; J. Webersfeld and E. Erb. J. Chem. Phys., 51, 328, 1954.

ASSOCIATION: Fiziko-tekhnicheskiy institut Kazanskogo filisla AN SSSR i Institut organicheskoy khimii AN SSSR Kazan' (Physico-

technical Institute of the Kasan' Branch of the AS USSR and

Institute of Organic Chemistry of the AS USSR Kazan')

SUBMITTED:

February 19, 1960

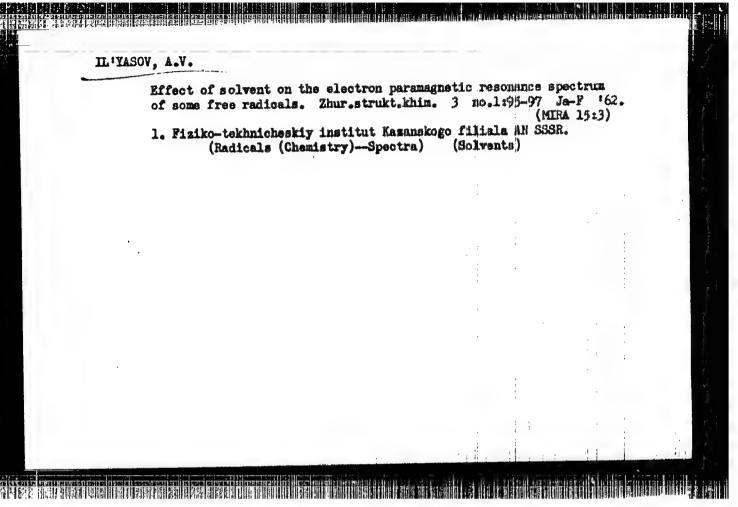
Card 4/6

ARBUZOV, B.A.; KATAYEVA, L.M.; KATAYEV, Ye.G.; II. YASOV, A.V.

Electron paramagnetic resonance studies of the dissociation of di-(2,4,6-triphenyl) phenyl paramale and di-(4,4,6-triphenyl) phenyldiselenide to free redicals. Inv. AN SSSE Otd, thin. nauk no.2:360-362 F '62.

i. Kazanskiy gosudarstvennyy universitet im. V.I. Ul'yanovalenina i Kazanskiy filial AN SSSE.

(Radicals(Chemistry))



s/020/62/144/003/027/030 B124/3101

AUTHORS:

Valitova, F. G., and Il'yasov, A. V.

TITLE:

The electron paramagnetic resonance in concentrated

α,α-diphenyl-β-picrylhydrazyl solutions

PERIODICAL:

Akademiya nauk SSSR. Doklady, v. 144, no. 3, 1962, 600-601

TEXT: The dependence of the relaxation time T_1 and T_2 on concentration was determined by continuous saturation at a frequency V= 460 Hc/sec in α, α -diphenyl- β -picrylhydrazyl solutions in benzene, toluene, and chloroform for concentration between 0.17 and 0.025 moles/liter at temperatures between 240 and 320°K. There is only a single paramagnetic absorption line with a distance of 4.3 oe between the inflection points which corresponds to the maximum concentration. The ratio $<\Delta {\rm H}^4>^{1/4}<\Delta {\rm H}^2>^{1/2}$ of 1.38 is indicative of a Lorenz-type absorption curve. When the concentration C is 0.025 moles/liter, exchange interactions become so small that the hyperfine structure characteristics reappear. T2 is calculated from the relation $T_2 = 1/\pi\sqrt{30}\nu$, where $\delta\nu$ is the line

Card 1/3

S/020/62/144/003/027/030 B124/B101

The electron paramagnetic ...

width in frequency units, whereas T1 is calculated from the saturation equation Z = $(1 + 0.25\gamma^2 H_0^2 T_1 T_2)^{-1}$, where Z is the saturation coefficient, γ is the gyromagnetic ratio, and H, is the amplitude of the high-frequency magnetic field. Relaxation time is found to be independent of the type of solvent used. The same order of magnitude of T, and T2 for concentrations of 0.17 moles/liter is indicative of a strong interaction exchange. T1 increases as compared to T2 in less concentrated solutions, and both become dependent on temperature. The heat-accumulator model developed by N. Bloembergen and S. Wang is used to interpret the results obtained. In the solution where the concentration is highest and the interaction exchange is large, the energy absorbed by the Zeeman system is transferred to the exchange system with the relaxation time T. F.T., where T, is the spin-lattice relaxation time and T_2 the spin-spin relaxation time. fact that the relaxation time is independent of temperature shows that it is not related to the Brownian motion of the paramagnetic melecules. On dilution, exchange is reduced and relaxation due to the Brownian motion of radical molecules increases. It is also found that the exchange frequency Card 2/3

The electron paramagnetic ...

B/020/62/144/003/027/030 B124/B101

 $\omega_{\rm a} \geqslant 10^{10}~{\rm sec}^{-1}$. There are 1 figure and 1 table. The most important English-language reference is: N. Bloembergen, S. Wang, Phys. Rev., 93,

ASSOCIATION:

Fiziko-tekhnicheskiy institut Kazanskogo filiala Akademii nauk SSSR (Physicotechnical Institute of the Kazan' Branch

of the Academy of Sciences USSR)

PRESENTED:

January 26, 1962, by A. Ye. Arbuzov, Academician

SUBMITTED:

January 24, 1962

Card 3/3

S/020/62/147/001/015/022 B106/B101

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Study of the free radical ...

(I). Data for the radical: yield 70-80%;

small bright-orange crystals with a melting point >180°C (decomposition); soluble in benzene, chloroform, alcohol, acetonitrile, glacial acetic acid and dioxane. In dilute solutions (< 10°) moles/l), the spectra show a hyperfine structure, the analysis of which proves that the unpaired electron in I remains mainly on the nitrogen atoms. A comparison of the e.p.r. spectrum of I with the spectrum of the α,α-diphenyl-β-picryl hydrazyl radical (DPPH) showed that the additional hyperfine structure is due solely to the protons of the α-phenyl groups. It may be explained by the interaction of the unpaired electron with the 2,4,6-protons of one of the two α-phenyl groups. The value obtained for the constant a of hyperfine coupling was 1.7 cersteds, and for ΔH_n 1.1 cersteds. The relative Card 2/4

Study of the free radical ...

\$/020/62/147/001/015/022 B106/B101

stability of related free radicals from the e.p.r. spectra are estimated by the method of J. A. Weil, K. V. Sane, J. M. Kinkade (J. Phys. Chem., 65, 710 (1961)) showed that I is chemically more stable than DPPH. Its stability may be due to steric factors reducing the possibility of chemical reactions with other substances. The values obtained from the e.p.r. spectra of I in finely crystalline state, which may contain solvent, were 15.7 ± 0.3 cersteds for ΔH at $295^{\circ}K$, 10.5 ± 0.3 cersteds at $77^{\circ}K$, 1.43° for r at $295^{\circ}K$, and 1.45 at $77^{\circ}C$ ($r = \langle \Delta H^4 \rangle 1/4 / \langle \Delta H^2 \rangle 1/2$). The g-tensor at $295^{\circ}K$ is: $g_1 = 2.0039 \pm 0.0001$, $g_2 = 2.0051 \pm 0.0001$, and $g_3 < g_1$. The considerable difference between these values and the g-factor of DPPH suggests that the molecular structure of the free radical considerably affects the residual spin - erbital coupling and anisotropy of the g-factor. There are 3 figures and 1 table. The most important English-language references are: M. N. Chen, K. V. Sane et al., J. Phys. Chem., 65, 713 (1961); B. Kubo, K. Tomita, J. Phys. Soc. Japan, 9, 888 (1954); F. K. Kneübuhl, J. Chem. Phys., 33, 1074 (1960).

Card 3/4

Study of the free radical ...

8/020/62/147/001/015/022 B106/B101

ASSOCIATION: Fiziko-tekhnicheskiy institut Kazanskogo filiala Akademii nauk SSSR (Physicotechnical Institute of the Kazan' Branch of the Academy of Sciences USSR); Khimicheskiy institut im. A. Ye. Arbuzova Akademii nauk SSSR (Chemical Institute imeni A. Ye. Arbuzov of the Academy of Sciences USSR)

SUBMITTED:

August 8, 1962

Card 4/4

CIA-RDP86-00513R000618520013-5" APPROVED FOR RELEASE: 04/03/2001

Hyperfine structure of ...

S/020/62/147/005/022/027

increases the electron density of the unpaired electron on the Nyatom.
Substitution of one methoxy group for one p-H atom of the *-phenyl group

CH_3-0-W-W- more probable than in a nonsubstituted radical. Substitution of NO2 for one p-H in the phenyl group

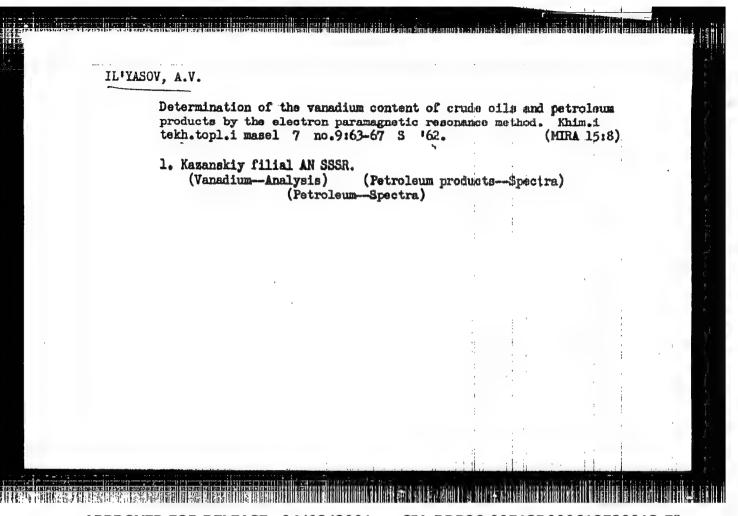
of triphenyl methyl causes polarization of the electron clouds of the

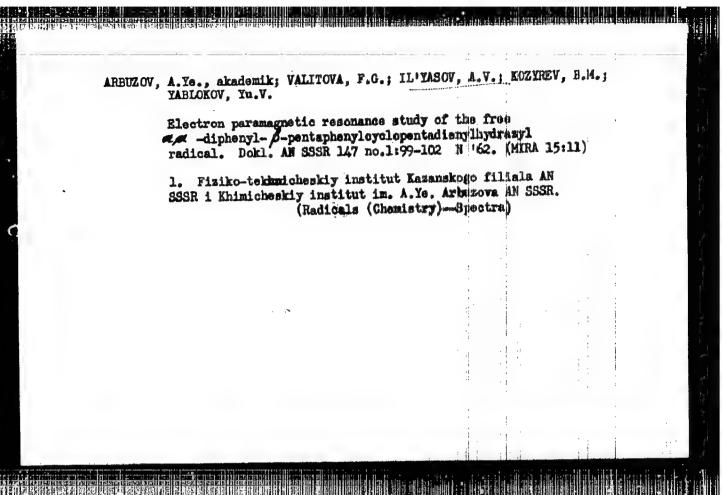
sequence: NA-C-, and NA-W- bonds. Polarization decreases in the following

sequence: NA-W- bonds. Polarization decreases in the following

of the unpaired electron on the Na atom revealed by the high A/W values,
The most important English-language references are: R. M. Denl, W. 3.

Koski, J. Chem. Phys., 31, 1138 (1959); N. W. Lord, S. W. Blinder, J. Chem.
Phys., 54, 1693 (1961); Y. Deguchi, J. Chem. Phys., 32, 1564 (1960).





IRRIHA, M.A.; IL'YASOV, A.Y.; MOZIREV, B.M.; MATEVOSTAH, R.O.;

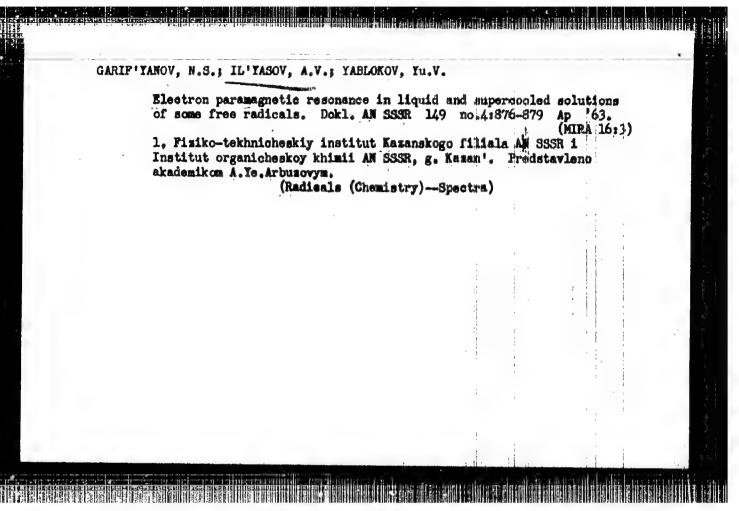
RYZIMANOV, Yn.M.: TABLOKOV, Yn.W.

Superfine structure of electron paramagnetic resonance spectsa of d.p. —diphenyl—pt-triphenylmethylhydramyl and its derivatices.

Dokl. AN SSSR 147 no.3:618-621 N '62.

1. Fixiko-tekimicheskiy institut kasanskogo filiala AN SSSR i Ural'skiy politekimicheskiy institut im. S.M. Kirova, Fredstavleno skademikom B.A. Arbusovym.

(Eydrasine) (Radicals (Chemistry)....Spectra)



KOZTREV, B.M.; YABLOKOV, Yu.V.; MATEVOSYAN, R.O.; IKRIA, M.A.;
IL'YASOV, A.V.; RYZHMANOV, Yu.M.; STASHKOV, L.I.; SHATRUKOV, L.F.

Electron paramagmetic resonance in substituted dipherylpicrylhydrasyls.
Opt. 1 spektr. 15 no.5:625-635 N '63. (KIRA 16:12)

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EPF(c)/FWP(1)/FWT(1)/EWT(m)/BDS-+AFFTC/ASD--PT-L/PG-4--PH/WW/JW/ 5/0020/63/150/001/0588/0591 10831-63 AP3000754 ACCESSION NR: Il yasov, A. V.; Garif yanov, N. S.: Timerov, H. Kh. AUTHOR: The nature of spin-lattice interaction in magmentically weak free radicals TITLE: Doklady, v. 150, no. 3, 1963, 588-591 AN SSSR. SOURCE: TOPIC TAGS: electron paramagnetic resonance, time of upin, leather relation, Alpha, Alpha-diphenyl-Beta-picryl-hydrazyl ABSTRACT: The electron paramagnetic resonance (e.p.r.) was intudied in solutions of free radicals of Alpha, Alpha-diphenyl-Beta-pieryl-hydlwzylland 2.2.6. tetranethylpentamethylene mitric oxidedin methanol, ethanol, henzene, kdiuene and mixtures of these in glycerin and in water. A study of solid (superconled) solutions (10 sup -2 to 10 sup -3 mol/1) indicated the times of spin lattice relaxation was independent of concentration and nature of solvents. The mechanism proposed by I. V. Aleksandrov and G. M. Zhidomirov (Zh. E. T. F., Wil. 147, 1961) provides for relaxation time in solid solutions of free radicals. Experiments run at ellevated temperatures indicated that collisions (brownian movement) in polar solvents (solvated radicals) were less effective on relaxation than in non-polar solvents (non-solvated radicals). Intensification of signal is not proportional to increase Card 1/2

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ACCESSION NR: AP3000754

in concentration of radicals, but much greater. This supports proposal by Mckennel (J. chem. phys. 25, 709, 1956) that isolated radicals have too long a releasablen time and are therefore saturated by small forces of the high frequency field and do not contribute to the e.p.r. signal. In these dilute collutions the mechanism is considerably dependent on the nature of the solvent. "The authors express thanks to B. M. Kozyrev for discussion of the results." Orig. art. has: 3 equations, 1 table, 1 figure.

ASSOCIATION: Fiziko-tekhnicheskiy institut Kazanskogo filiala fikademii pauk ISSR (Physical-Technical Institute of the Kazan Bramch of the Academy of Sciences ISSR). Institut organicheskoy khimii Akademii nauk SSSR Kazan (Imstitute of Organic Chemistry, Academy of Sciences SSSR)

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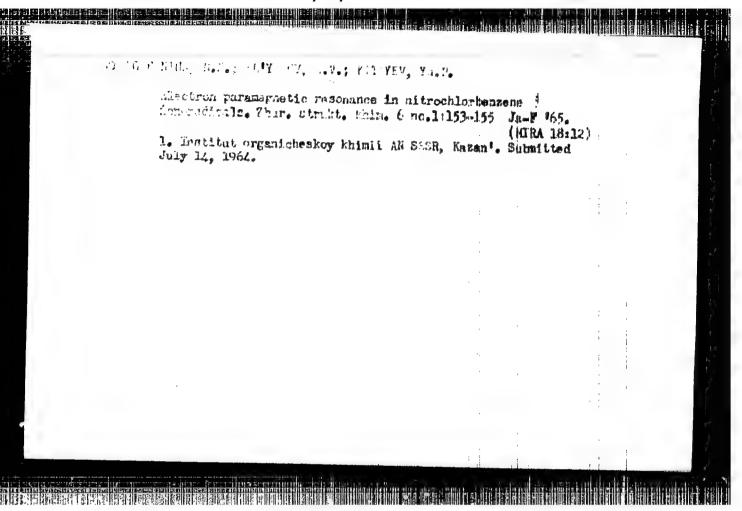
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VOZDVIZHENSKIY, G.S.; GUDIN, N.V.; SHAPHIK, M.S.; GAHIF'YANOV, N.S.;
IL'YASOV, A.V.

Electron paramagnetic resonance study of the electrode processes of copper complexes with organic amino derivativas. Zhur. fis. khim. 38 no.6:1682-1685 Je '64. (MIRA 18:3)

1. Kazanskiy khimiko-tekhnologicheskiy institut imeni Kirova i Institut organicheskoy khimii AN SSSR, Kazan'.



VALITOVA. F.G.; ILIVASOV, A.V.; SOTNIKOVA, N.N.; BARCH. DINA, S.Yu.

Electron paramagnetic resonance study of electrochemically generated radicals of some hydrasines. Shur.strukt.khim.
6 no.52777-779 S-0 '65.

1. Institut organicheskoy i fizicheskoy khimil AN SSSR, Kazan'.

VOZDVIZHENSKIY, G.S.; GUDIN, N.V.; SHAPNIK, M.S.; IL'YASOV, A.V.; GARIF'YANOV, N.S. (Kazan')

Klectron paramagnetic resonance study of electrode processes in aqueous solutions of copper complexes. Zhur. fiz. khim. 39 no. 1: 64-67 Ja *65 (MIRM 19:1)

1. Institut organicheskoy khimii AN SSSR, Kazan'. Submitted January 10, 1964.

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AUTHOR: Illyasov, A. V.: Levin V.
AUTHOR: Il'yasov, A. V.; Levin, Ya. A.; Sotnikova, N. N.; Valitova, P. C. 85
ORG: Institute of Organic and Physical Chartes
ORG: Institute of Organic and Physical Chemistry, AN SSSR, Karan' (Institut 72
TITIE: Electrochemical generation of hydrazyl radicals
SOUTHORN
SOURCE: Teoreticheskaya i eksperimental naya khimiya, v. 2, no. 1, 1966, 142-143
TOPIC TAGS: 01004 1, 1966, 142-143
TOPIC TAGS: electrochemistry, free radical, hydrazine derivative, electrolytic cell, resonator
resonator paramagnetic resonance, redor resetter
resonator/RE-1301
ABSTRACT: The American
β-picrylhydragyl (DPPH) are obtained by treating the corresponding hydragines. bility of abtainst or other oxidizing agents. The authors statistics
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electrolyte. To improve the recollection used as the supporting
of oldered parameter and the same the
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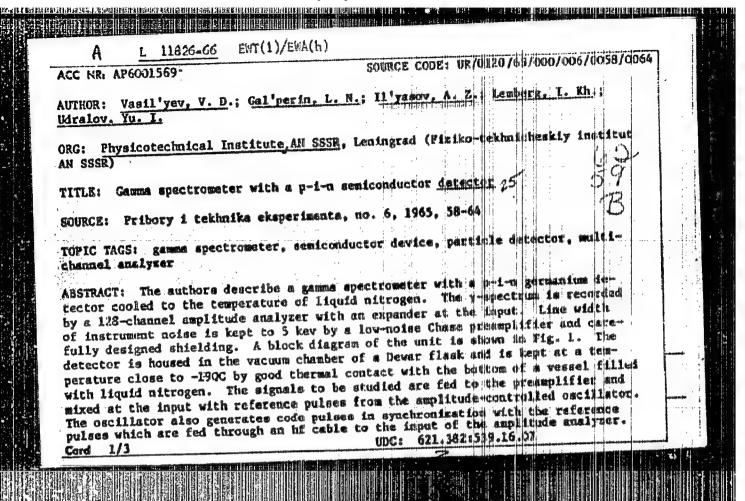
solutions were degassed by the freezing method. The formation of hydraxyls in electrochemical oxidation of the original compounds can be depicted by the scheme:

$$Ph_2N - NH - R + OH^- \stackrel{?}{=} Ph_2N - N - R + H_2O$$
,
 $Ph_2N - N - R \stackrel{?}{=} Ph_2N - N - R + O$.

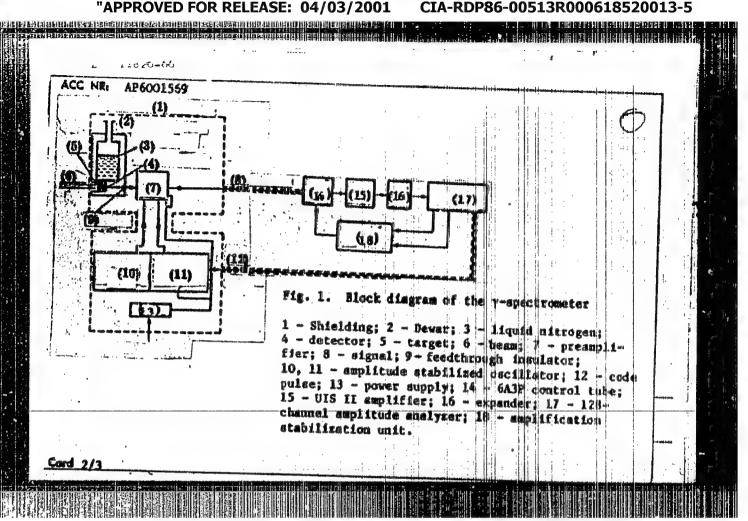
Thus, the authors have shown that electrochemical oxidation as well as electrochemical reduction of compounds of the diphenylpicrylhydrasine type lead to the formation of free radicals, the properties and structure of which can be studied by the electron paramagnetic resonance method, TPRS

SUB CODE: 07 / SUBM DATE: 21Jun65 / CRIC REF: 006 / OTH REF: 004

Card 2/2 /11/2



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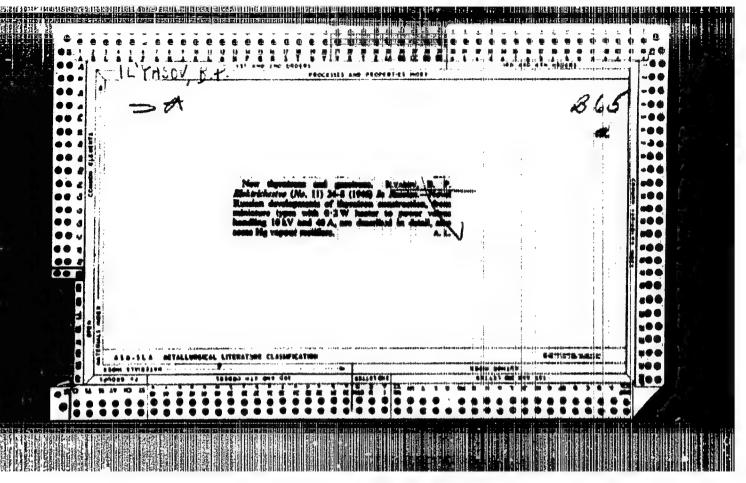


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ACC NR: AP6001569 The code pulses separate the reference pulses from the detector starists after amplification. These same code pulses prevent registration of the reference pulses when the detector signals are being recorded. Pulses from a second amplitudecontrolled oscillator may also be fed to the preamplifier input for simulating detector signals when checking the operation of the device. From the output of the preamplifier, the signals being studied and the reference pulses are fed to the to the third grid of a 6A3P tube, which controls amplification during stabilization. Amplification control voltage from the stabilisation unit is find to the first grid of this tube. The signals are then amplified by a UIS-II amplifier and fed through the expander to the amplitude analyzer. The various meetions of the unit are described in detail, with diagrams of the cooling unit, low-noise preampliffer, expander, stabilization circuit, and output stage of the amplitude controlled oscillator. Tests showed that continuous-duty stability of the analyzer is better than 0.15% with no apparent effects of interference from the cyclotron with which it is designed to be used. The authors thank S. M. Ryvkin, O. A. Matveyev, and N. B. Stroken for graciously supplying experimental detector models. Orig. art. has: 8 figures. [01] SUB CODE: 40,09/SUBM DATE: 170ct64/ ORIG REF: 003/ OTH REE: DOL ATD PRESS:9//

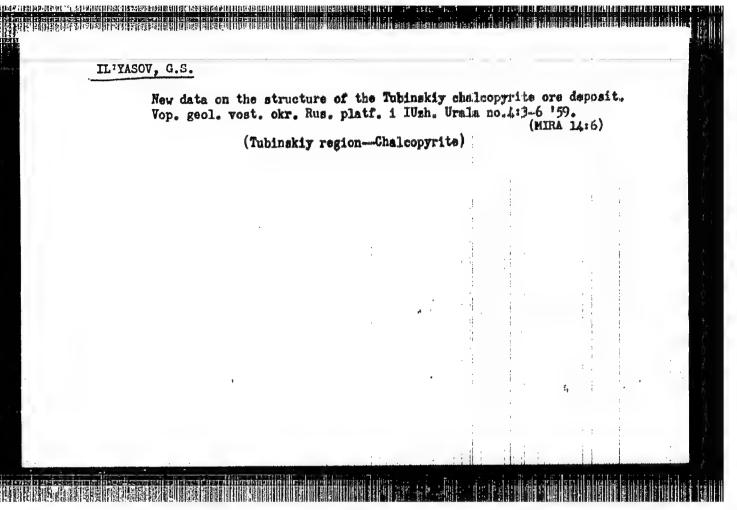
__IL'YASOV, B., inzh. (g. Ashkhabad)

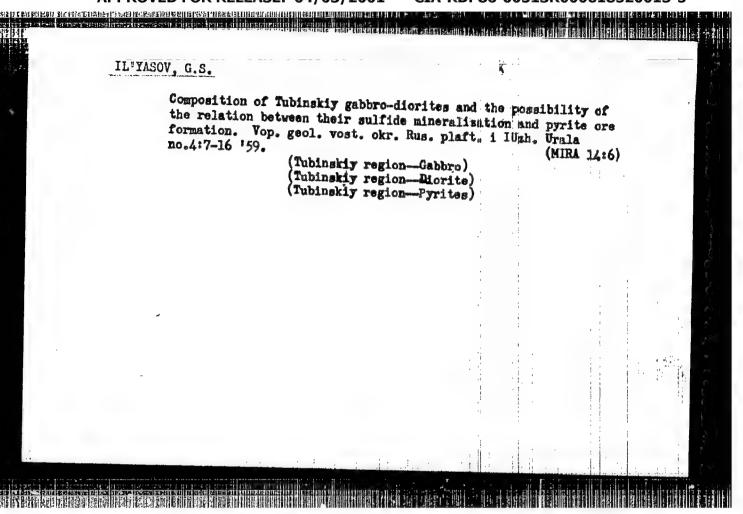
Seismic resistance of sectional irrigation structures.
Gidr. i mel. 17 no.7:15-25 Jl '65. (MIPA 18:12)

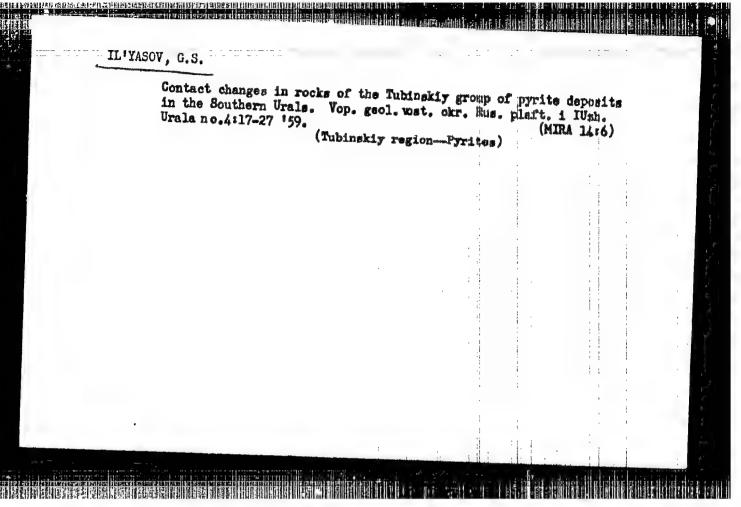


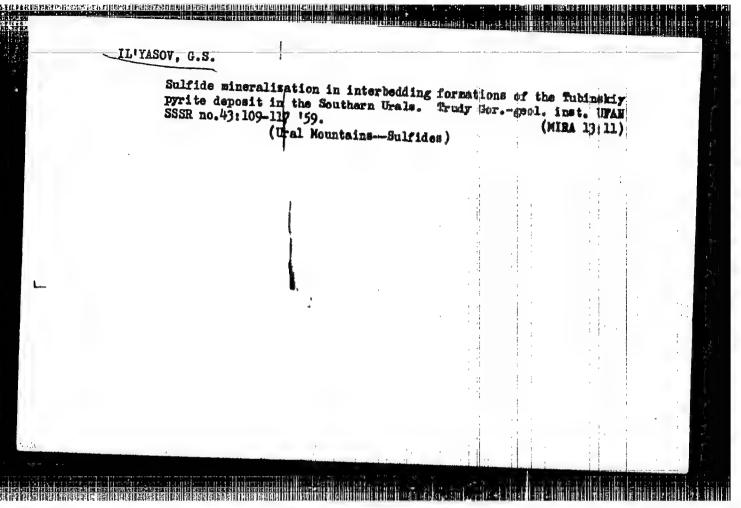
IL:YASOV, G.S., Cand Geol Min Sci — (diss) "Geological structure and age of the Tubinskiy pyrite deposit in the southern Urals." Ufa, 1959, lh pp (Acad Sci USSR. Rak Ural Affiliate) 150 copies (KL, 35-59, 113)

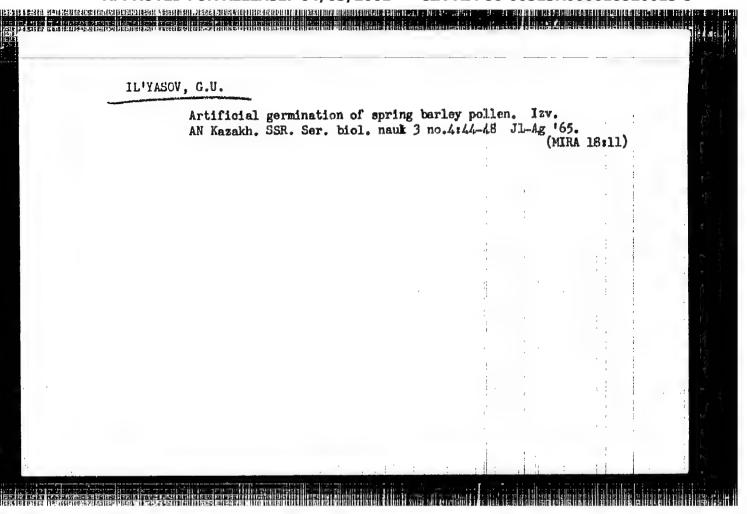
- 21 -

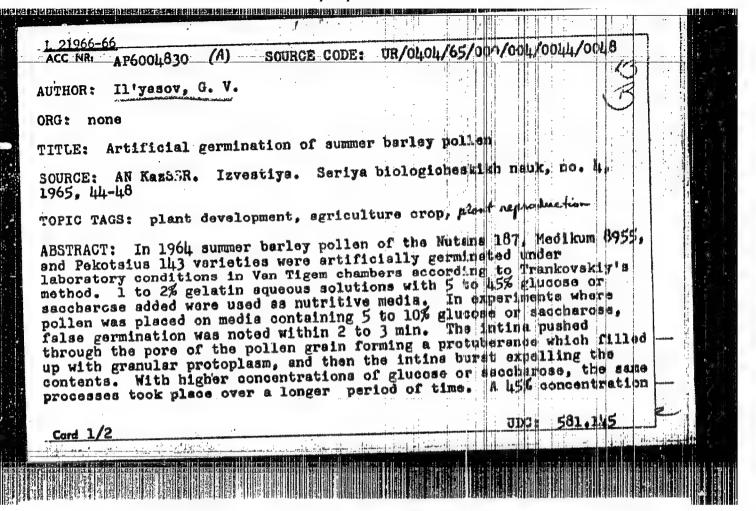








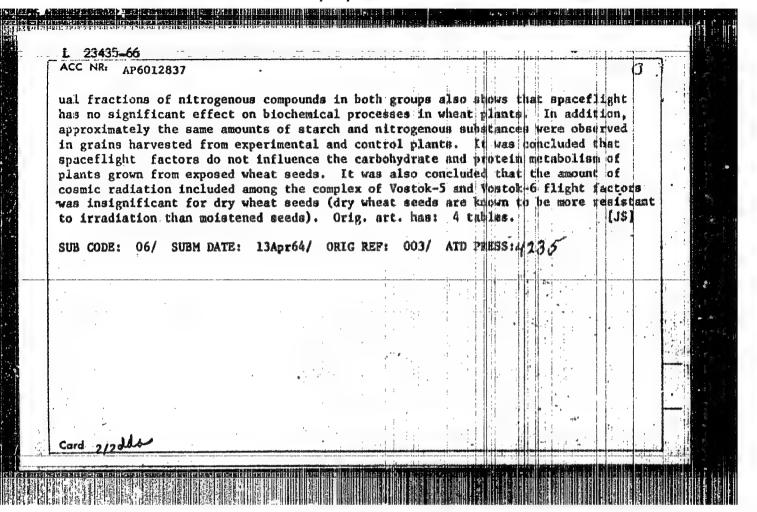




1. 21966-66 ACC NR: AP6004830 caused plasmolysis of the pollen grains. The experiments were repeated over 8 days with the same results. Another series of experiments was conducted according to methods described by S. Antony and H. V. Herlan (1920) in which the pollen was placed on the under side of a microscope slide cover instead of the glass slide and depression slides covered with large cover slides (32 x 40 mm) were used instead of Vin Tigem chambers. The slides were placed on a window sill and observed 3, 5, and 10 min following exposure. After 5 min most of the visble pollen grains germinated and the rest remained unchanged throughout the experiment. Findings show that barley pollen is sensitive to insufficient moisture as well as excessive moisture. Under favorable conditions barley pollen germinates within the first initials. Exposure of pollen grains to dry air leads to change in form; however, with immediate transfer of the pollen to a moisture plamber, viability of pollen is preserved and it germinates normally. With more prolonged exposure to dry air, the pollen loses its germination capacity. S. Antony and H. V. Harlan's method is recommended for ertificial germination of barley pollen. Orig. art. hast 4 fightes. OTH REP SUB CODE: 06/ SUBM DATE: none/ ORIG REF: 001/ 001 Card 2/2 ULP

ika (1996) Penung dengangangkangangan Per'anni aksala, sebuluh menghini dalah menghinin dan 1996. Per

23435-66 FSS-2/EnT(1)/EEC(k)-2/EMA(d) SCTB ACC NR: AP6012837 SOURCE CODE: UR/029/1/66/004/002/0320/0320 AUTHOR: Il'ina, G. V.; Kuznetsova, N. N.; Rydkiy, S. C.; Vyuutskily, Y. G. ORG: none TITLE: The effect of spaceflight factors on wheat seeds and plants grown from them SOURCE: Kosmicheskiye issledovaniya, v. 4, no. 2, 1966, 320 323 TOPIC TAGS: space biology, radiation effect, germination, whitat, carbohydrate metabolism, protein metabolism, plant physiology ABSTRACT: A study was made of the growth and development of wheat plants grown from seeds exposed to spaceflight factors on the Vostok-5 and Vos ok-6 lights. Experimental and control batches of wheat seeds ("Krasnozerna" var dty) were cultivated in fertilized soil under controlled humidity conditions. The energy of germination of seeds was determined and biochemical analysis was made of the plants in the following growth phases: seedling stage, tillering stage, and late flowering stage. A slight tendency to depression of germination was observed in experimental seeds (10% fewer sprouts on the first day of counting). Study of plant growth and accumulation of dry mass showed no difference between experimental and control groups. Biochemical analysis of plants showed insignificant variations in the content of individual sugar fractions, and analogous changes in soluble carbohydrate content in both experimental and control seedlings. The similarity of changes in nitrogen content and in individ-581.057

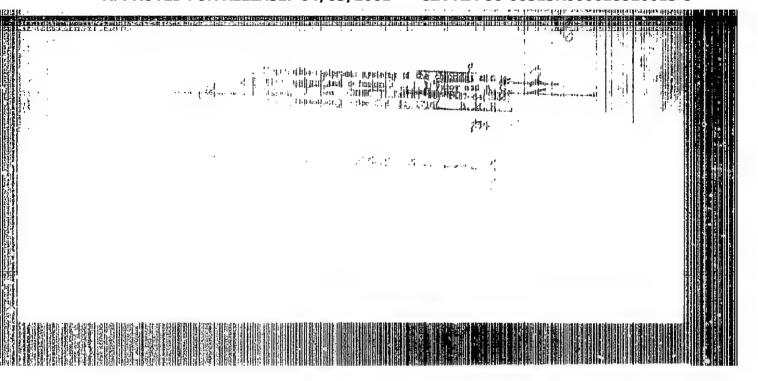


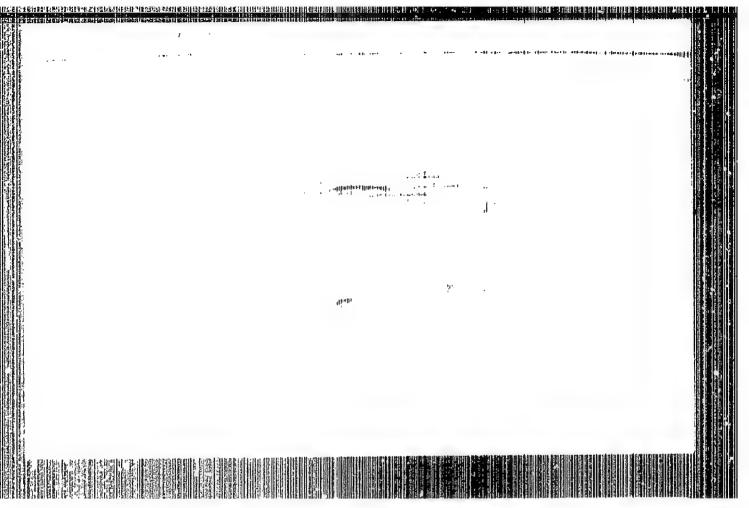
IL'YASOV, I.I. ---

"Physicochemical Analysis of Choride-Iodide Exchange in Fusions of Salts of Mono- and Divalent Metals." Cand Chem Sci, Novocherkassk Polytechnic Inst, Novocherkassk, 1954. (RZhKhim, No 20, Oct 54)

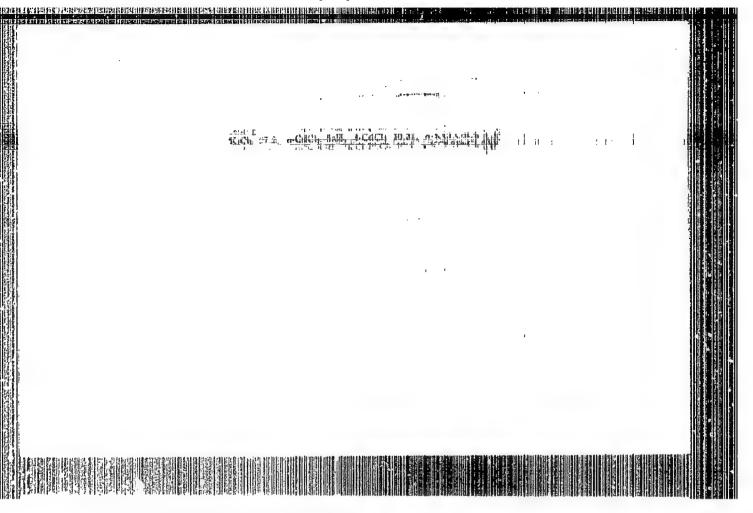
Survey of Scientific and Technical Dissertations Defended at UBSR Higher Educational Institutions (10)

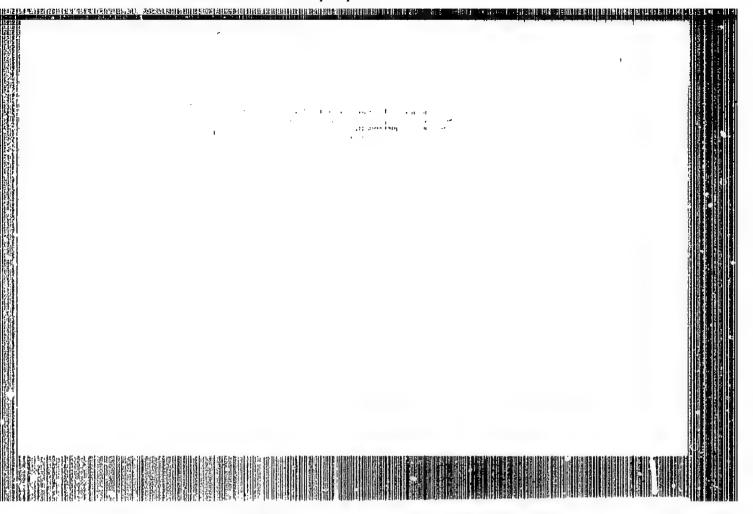
SG: Sum. No. 481 5 May \$5





APPROVED FOR RELEASE: 04/03/2001 CIA-RDP86-00513R000618520013-5"





APPROVED FOR RELEASE: 04/03/2001 CIA-RDP86-00513R000618520013-5"

"The Surface of Crystallization in the Constitutional Diagram of the Ternary System Composed of the Chlorides of Schium, Potassium, and Cadmium, by I. I. Il'yasov, A. K. Bostandriyan, and A. G. Bergman, Rostov-na-bonu Engineering-Construction Institute, Zhurnal Heorganicheskoy Khimii, Vol 2, No 1, Jan 57, pp 172-178

The ternary system Na, K, Cd/Cl was subjected to investigation. The constitutional diagram which was obtained differed in some essential reconstitutional diagram which was obtained differed in some essential respects from that determined by non-USSR scientists. It was established spects from that determined by non-USSR scientists. It was established that the stable compound KCl-CdCl2 is formed, which melts without decomposition, and that the unstable compounds hall (CdCl2 and MaCl-CdCl2) with melt with decomposition, are also formed.

I LYASOV, [.].
Bergman, A.G. and Iliasov, I.I. 575 Fusion Diagram for the Reciprocal System of Cadmium and Potassium Chlorides and Iodides. (Diagramma Playkosti Vzaimnoy Sistemy iz Khloridov i Yodidov Kadmiya i Kaliya.) AUTHORS: "Zhornal Neorganicheskoy Kwimii" (Journal of Inorganic Chemistry TITLE: Vol.11, No.2, pp.395-406. (U.B.S.R.) (95) The system K, Cd | Cl, I has clearly developed complex forma-PERIODICAL: tion of the binary-system components and polymorphism, The crystallisation surface has a fairly complicated relief and consists of ten fields, meeting in six mon-variant points. ABSTRACT: Because of the decomposition of CdI2 the investigation of the system reported was restricted to temperatures below 550°C. Carbon dioxide was continuously passed into the melt to minimise Study of the liquidus diagram of the Cill2 - K2I2system showed the existence of the compound CdI2.2KI and CdI2.KI, melting the existence of the compound CdI2.2KI and CdI2. KI, melting with decomposition at 223 and 272°C, respectively. The system with decomposition at 223 and 272°C, respectively. The system cdCl2 - CdI2 has a cutectic at 360°C and 31% CdCl2 and a homeomorphous transformation for CdCl2 at 460°C. In the system cdCl2 - KaCl2 the compound CdCl2 4KCl was found, melting at morphous transformation for CdCl2 at 40000, in this bystam CdCl2 - K2Cl2 the compound CdCl2.4KCl was found, melting at 428°C without decomposition. Eleven different fields of crystallisation were found in the reciprocal system K, Cd | Cl, I card 1/2 Territoria de la contribidad de distribilitativa de distribuciones de Carl of the party map we have interested

Fusion Diagram for the Reciprocal System of Cadmium and Potassium Chlorides and Iodides. (Cont.)

The complex CdCl₂.KCl (α , β) is stable and occupies an area of 14.72%. The compounds α and β CdCl₂.4KCl, CdI₂.2KI and CdI₂.KI, which melt with decomposition, also retain their stability inside the system and have a common crystallisation curve with the complex CdCl₂.KCl and component CdI₂. In the system complex-formation predominates over exchange reaction. The system is divided into four phase triangles by the three adiagonal triangulating sections: CdCl₂.KCl - CdI₂.CdCl₂.KCl - CdI₂.2KI and CdI₂.2KI - K₂CI₂. There are eight references, four of them Russian.

10 Figures, 7 Tables.

The work was carried out at the Engineering-Construction Institute, Rostov on Don.

Received 8 May, 1956.

Card 2/2

Category: USSR / Physical Chemistry

Thermodynamics. Thermochemistry. Equilibrium. Physico-

chemical analysis. Phase transitions.

B-8

Abs Jour: Referat Zhur-Khimiya, No 9, 1957, 29946

Author : Il'yasov I. I., Berguan A. G.

Inst : not given

1 L 41 - - F

Title : Irreversibly-Mitual System of Chlorides and Iodides of Sodium and

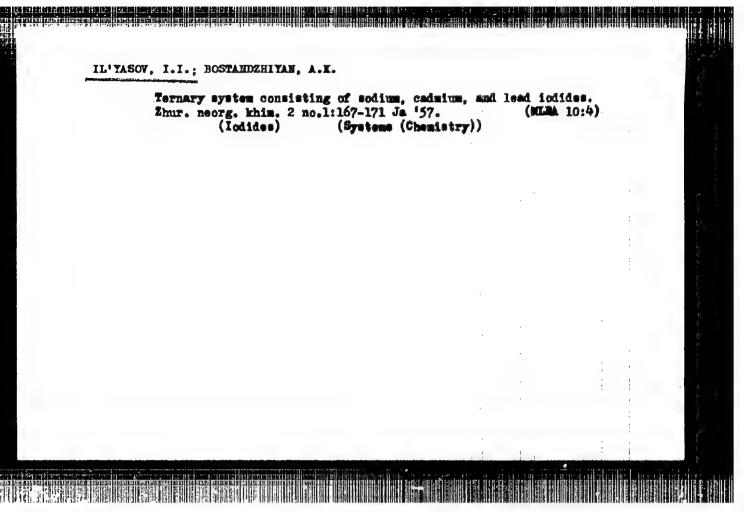
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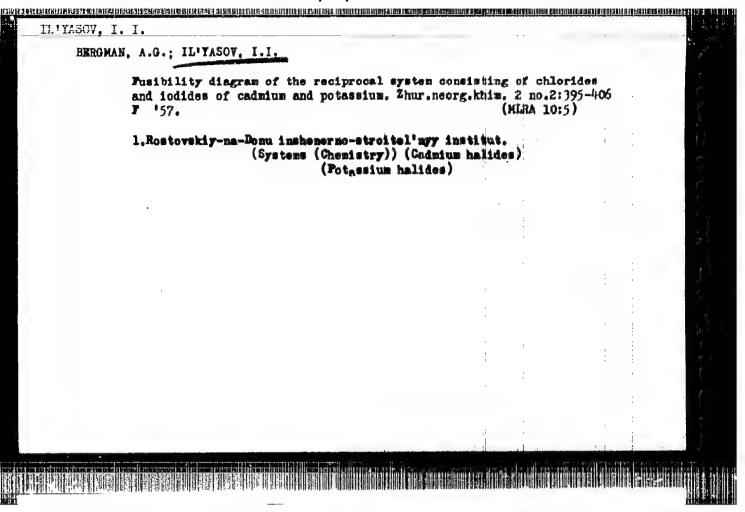
Orig Pub: Zh. obshch. khimii, 1956, 26, No 5, 1268-1296

Abstract: Study of the mutual system Na, Cd / Cl, I. The stable diagonal section is Na₂Cl₂- CdI₃; the subordinate, adiagonal, triangulating secant is Na₂Cl₂- CdI₃: 2NaI. Positive conditional thermal effect of reaction, equal to 6.7 kcal/equivalent, indicates the irreversible nature of exchange reaction in the system. There is confirmed the formation of the previously ascertained compound CdCl₂: 2NaCl₃, which melts with decomposition; transition point at 433° and 37.5% Na₂Cl₂. There was ascertained a compound CdI₂: 2NaI₃, melting with

Card : 1/2

-63-





CIA-RDP86-00513R000618520013-5 "APPROVED FOR RELEASE: 04/03/2001

Il- YASOV,

USSR/Physical Chemistry. Thermodynamics, Thermochemistry, B-8 Equilibria, Physical-Chemical Analysis, Phase Transitions.

Abs Jour: Ref Zhur-Khimiya, No 5, 1957, 14696

I. I. Il' yasov, A. G. Bergman Author

Inst

Reciprocal System of Potassium and Lead Chlorides and Title

Iodides with Interior Heterocomplex

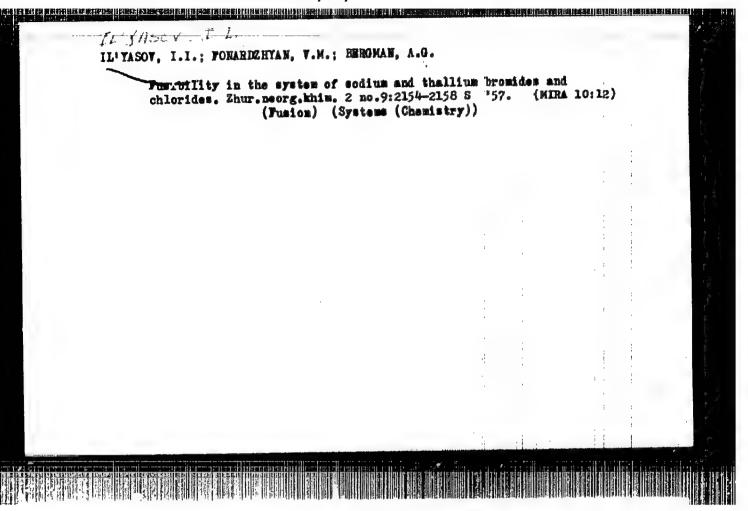
Zh. obshch. khimii, 1956, 26, No 4, 981-991 Orig Pub:

The fusibility graph of the reciprocal system E, Pb // Cl, I (I) was studied. A complex in the form of a Abstract: binary compound (surmised composition PbI2 · KC1) was detected on the stable diagonal KC1-PbI2 of the system I. The presence of the complex PbCl2 PbI2 fusing without decomposition was established in the system PbCl2-PbI2. The system I is divided into 8 phase triangles by the triangulating diagonal section PbI2-K2Cl2 and 6 adiagonal secants. The liquidus area consists of 9 fields and one

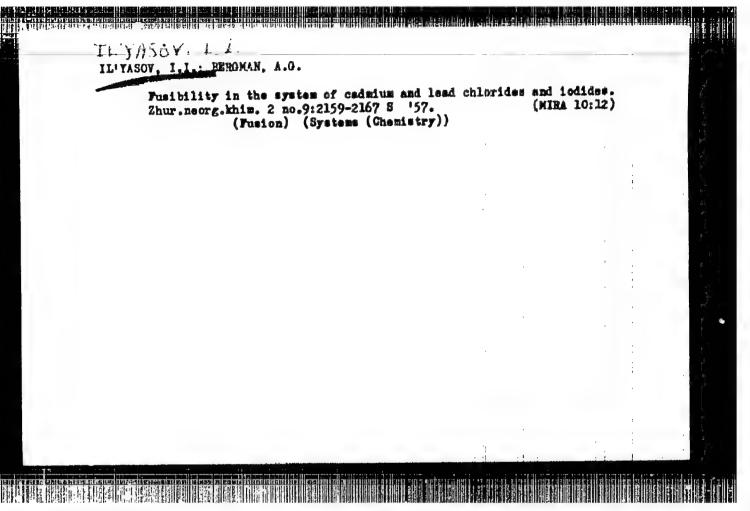
additional field dependent on the presence of PbI2

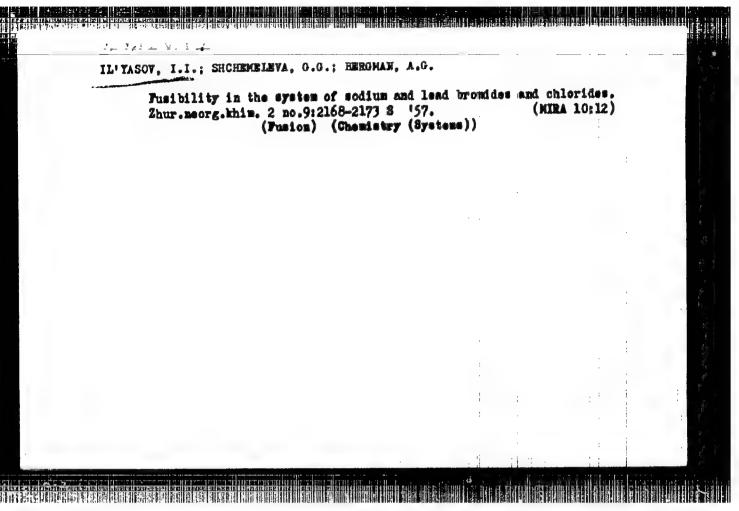
Card 1/2

Card 2/2



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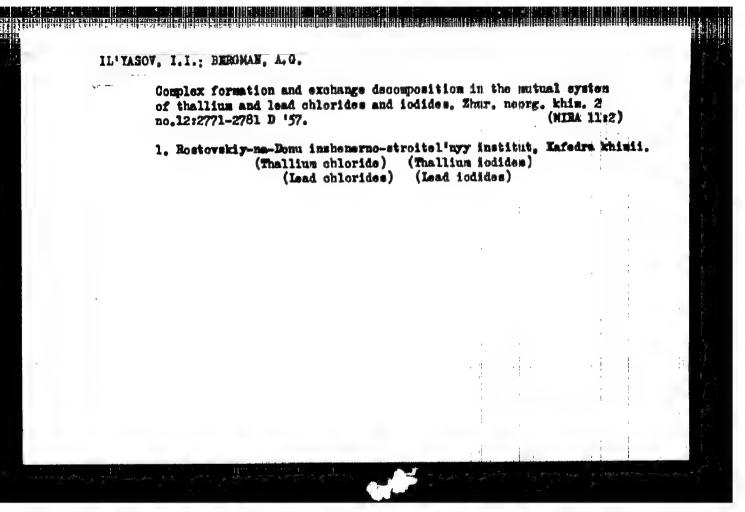


IL'YASOV, I.I.; ROZHKOVSKAYA, L.V.; RERGMAN, A.G.

Fusibility in the ternary sutual system of cadmium and lead chlorides and browides. Zhur.neorg.khim. 2 no.9;2174-2177
S '57. (MIRA 10:12)

l.Rostovskiy-na-Domu Inshenerno-stroitel'nyy institut.

(Fusion) (Systems (Chemistry))



SOV/78-4-4-33/44 5(4) Il'yasov, I. I., Shohemeleva, G. G., Bergman, A. G. AUTHORS: The Behavior of the Ternary System of Sodium, Cadmium and Lead Bromides in the Melting Process (Playkost' traynoy sistemy iz TITLE: bromidov natriya, kadmiya i svintsa) Zhurnal neorganicheskoy khimii, 1959, Wol 4, Wr 4, pp 906-908 PERIODICAL: (USSR) The system Na, Cd, Pb || Br was investigated by a visual polythermal method. The binary systems Naghra-FbBra, NagBra-GdBra ABSTRACT: and CdBr2-PbBr2 were checked and completed. Six intermal sections of the ternary system were injestigated; the results are contained in figure 1 and table 2. The melting diagram of this system consists of three main crystallization ranges. A range with $\alpha-$ and $\beta-$ homeomorphous differences appears within the range of Na Bro. In the system Na Bro-PbBro a sutectic cocurs at 324° with 9.7% Na Br 2. The translition point of the α - and β -homeomorphous form is located at 380° with 17% Nu₂Br₂. The system CdBr - PbBr forms a sutschic at 340° with Card 1/2

SOV/78-4-4-33/44
The Behavior of the Ternary System of Sodium, Cadmium and Lead Broatdes in the Melting Process

15% CiBr2. The melting points within the binary systems

PhBr - Na Br and PbBr - CdBr are given in a table.

There are 2 figures, 2 tables, and 8 references, 7 of which

are Sowiet.

SUBMITTED:

December 26, 1957

Card 2/2

SOV/78-4-4-34/44 Il'yasov, I. I., Mirsoyapov, V. N., Korbtkov, Yu. V. 5(4) AUTHORS: The Ternary System of Sodium, Potassium and Cadmium Bromid+8 (Troynaya sistema 12 bromidor natriya, kaliya i kadmiya) TITLE: Zhurnal neorganicheskoy khimii, 1959, Vol 4, Nr 4, pp 909-9:2 PERIODICAL: (USSR) The system Na, K, Cd || Br was investigated by a visual polythermal method. The binary systems Ma2Br2-K2Br2, Ma2Br2-CdBr2 ABSTRACT: and K2Br2-CdBr2 were oneoked, and it was found that in the system K2Br2-CdBr2 there is only one dompound with the composition KBr. 2CdBr2. This compound mells incongruently at 360°. Seven internal sections of the ternary system were investigated. The crystallization surface of this system consists of the ranges CdBr2, KBr.CdBr2 and the solid solutions [Na, K]Br, which decompose inside the system above 550°. The internal sections and the melting diagram of the system Ma, K, Cd | Br are given in figures 2 and 3. The melting points within the Card 1/2

SOV/78-4-4-34/44

The Ternary System of Sodium, Potassium and Cadmium Bromides

system CdBr,-K,Br, are contained in a hable; the seven internal

sections under investigation are characterized in a further table. There are 3 figures, 2 tables, and 6 references, 5 of

which are Soviet.

SUBMITTED:

December 30, 1957

Card 2/2

#07/70-4-4-35/44 5(4) Il'yasov, I. I., Bergman, A. G. AUTHORS: Complex Formation in the Reciprocal System of Chlorides and Iodides of Cadmium and Thallium (Kompleksoobrazovaniye po TITLE: vzaimnoy sisteme iz khloridov i yodidov kadmiya i talliya) Zhurnal neorganicheskoy khimii, 1959, Vd2 4, Mr 4, pp 913-919 PERIODICAL: (USSR) The reciprocal system T1, Cd | C1, J was investigated by a visual polythermal method. Pirst, the binary systems CdG12-CdJ2, ABSTRACT: CdJ2-Tl2J2. C4Cl2-Tl2Cl2 and Tl2J2-Tl2Cl2 ware investigated. In the system OdCl2-Tl2Cl2 the compound OdCl2.TlCl with the melting point 430° is formed. The unstable diagonal sections Ti2Cl2-CdJ2 and CdCl2-Ti2J2 were investigated; the results are given in figure 2. The triangulating nom liagonal sections from the top of the complex CdCl2.TlCl are given in figure 3. The section $CiCl_2$. TiCl-Tl₂J₂ consists of three branches; α - and β -OdCl₂-TlCl and Tl₂J₂- The section CdCl₂-TlCl-OdJ₂-4TlJ is Card 1/3

SCV/78-4-4-35/44

Complex Formation in the Reciprocal System of Chlorides and Indides of Cadmium and Thallium

characterized by polymorphous transformation of the branch CdCl₂.TlCl at 372° and 21.5% (CdJ₂.4TlJ). The following branches were found in the section CdCl₂.TlCl-CdJ₂: α-CdCl₂.TlCl, β-CdCl₂.TlCl, CdCl₂ and CdJ₂, which inversent at 372° and 21.5% CdJ₂, 330° and 35% CdJ₂, and 320° and 69% CdJ₂. Apart from the diagonal and triangulating sections thilteen internal sections were investigated, the melting diagrams of which are given in figures 4, 5 and 6. The crystallization surface of the system covers six physiallization ranges. The nature of the melting diagram shows that complex formation provails in the system Tl, Cd ||Cl, J. A characterization of the binary system CdJ₂-Tl₂J₂ and of the diagonal sections Tl₂Cl₂-CdJ₂ and CdCl₂-Tl₂J₂ by the melting points is given in a table; the three eutectic points and the point of transition of the system Cd, Tl ||Cl, J are contained in another table. There are 6 figures, 2 tables, and 12 references, 8 of which are Soviet.

Card 2/5

507/78-4-4-35/44 Complex Formation in the Reciprocal System of Chlorides and Iodides of Cadmium and Thallium

ASSOCIATION: Rostovskiy-na-Donu inzhenerno-stroitel'nyy institut

(Restov-na-Donu Institute of Construction Engineering)

SUBMITTED: January 15, 1959

Card 3/3

CIA-RDP86-00513R000618520013-5

SOV/76-4-9-25/44

AUTHORS: Bostandzhiyan, A. K., Il'yasov, I. I., Bergman, A. G.

TITLE: The Fusibility in a System of Chlorides and Bromides of Potassium and Lead

PERIODICAL: Zhurnal neorganicheskoy khimii, 1959, Vol. 4, Dr. 9, pp 2079-20 (USSR)

ABSTRACT: Before the combined system mentioned in the title is dealt with the melting curves of the binary systems K₂Cl₂ - PbCl₂

K₂Br₂ - PbBr₂ (in accordance with the data given by S. D. Gromakov, reference 2) F. Cl. - K. Response to the content of the data given by S. D. Gromakov,

reference 2), $K_2Cl_2 - K_2Br_2$ and $PbCl_2 = FbBr_2$ (in contrast with the data given by L. I. Favorskiy, reference i) are given in figure 1. In the combined system two diagonal and four interior sections were investigated (Table 1, Figs 2-4). In the four crystallization ranges K [Cl,Br], 2K [Cl,Br]. Fb [Cl,Br], K [Cl,Br]. 2Pb [Cl,Br]_2 and Pb [Cl,Br]_2 are formed. The system under examination belongs to the group of mutual systems with

Card 1/2

The Fusibility in a System of Chlorides and Bromides of Potassium and Lead

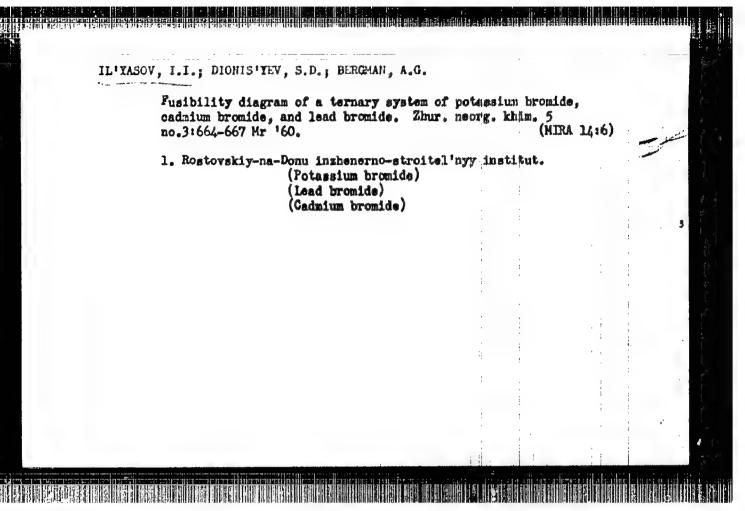
SOT/78-4-9-25/44

complex formations of the belt type in which all components and compounds of the sides opposite one another form stable continuous solid solutions with each other. There are 4 figures, 1 table, and 9 references, 7 of which are Soviet.

SUBMITTED: April 30, 1958

Card 2/2

5(2) 507/76-4-9-26/44 Il'yasov, I. I., Bergman, A. G. AUTHORS: The Fueibility in the Ternary System of Iodides of Sodium, FITLE Potassium, and Lead Zhurnal neorganicheskoy khimii, 1959, Vol 4, Nr 9, pp 2083-200 PERIODICAL: (USSR) After a short characterization of the binary systems Nagula - Kadas ADSTRACT: Na2J2 - PbJ2, and K2J2 - PbJ2 (Fig 1) a report is made on the investigation of the ternary system mentioned in the title (Fig 2, Tables 1, 2). Within this system an Enterior field can be clearly distinguished which borders on all the other components and the double compound EJPbJ2, melts incongruently, and has approximately the following composition: KJ.2NaJ.2FbJ, . It was found that the solid solutions of NaJ and KJ within this system decompose already below 5000. The four nonvariant points are given in table 3. There are 4 figures, 3 tables, and 7 Soviet references. SUBMITTED: April 30, 1958 Card 1/1



8/078/60/005/05/24/057 8004/8016

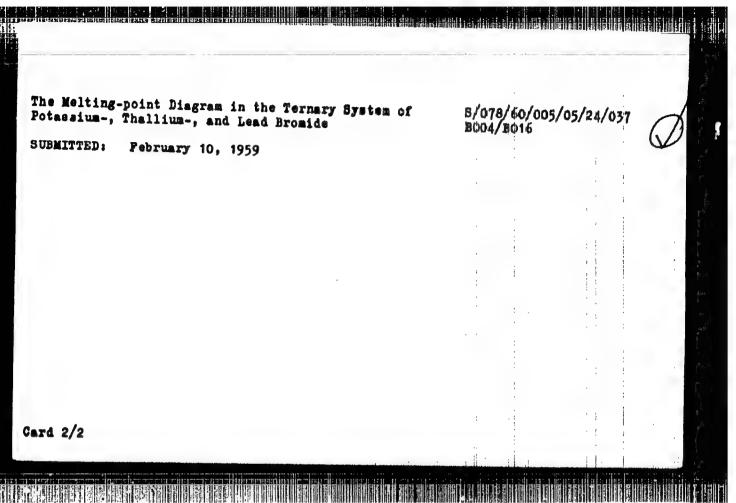
AUTHORS: Dionis'yev, S. D., Il'yasov, I. I., Bergman, A. G.

TITLE: The Melting-point Diagram in the Ternary System of Potassium-, Thallium-, and Lead Browide

PERIODICAL: Zhurnal neorganicheskoy khimii, 1960, Vol. 5, No. 5, pp. 1135 - 1138

TEXT: After giving a short survey of the binary systems $K_2Br_2 + Tl_2Br_2$; $K_2Br_2 - PbBr_2$, and $Tl_2Br_2 - PbBr_2$, and referring to the papers by A.P.Rostkovskiy (Ref. 2), and L. I. Favorskiy (Ref. 5), the authors report on their investigation of 14 sections (Tables 1,2, Fig. 1) of the ternary system. The resultant melting-point diagram is shown in Fig. 2, the crystallization scheme in Fig. 3. The melting-point diagram has a complicated structure owing to the formation of limited solid solutions between KBr and TlBr and stable, continuous, solid solutions between TlBr.2PbBr₂ and EBr.2PbBr₂. The phase diagram is divided into 5 phase triangles with 3 invariant points (Table 3). There are 3 figures, 3 tables, and 6 references, 5 of which are Soviet.

Card 1/2



IL'YASOV, I.I.; SHCHERLIYA, G.O.; REIGHAN, A.G.

Fusibility diagram of a ternary mystem consisting of sodium, potassium, and lead bromides. Ehur.neorg.khim. 5 no.6:1254-1256 [MIRA 13:7]

1. Rostovskiy-na-Donu inshenerno-stroitel'nyy institut. (Sodium bromide) (Potassium bromide) (Lead bromide)

77851 5.4110

Il'yasov, I. I., Bergman, A. AUTHORS:

Physico-Chemical Analysis of Systems Containing Salts TITLE:

of Organic Acids. I. Mutual System Consisting of

Chlorides and Acetates of Sodium and Potassium

Zhurnal obshchey khimii, 1960, Vol 30, Nr 2, pp 355-358 (USSR) PERIODICAL:

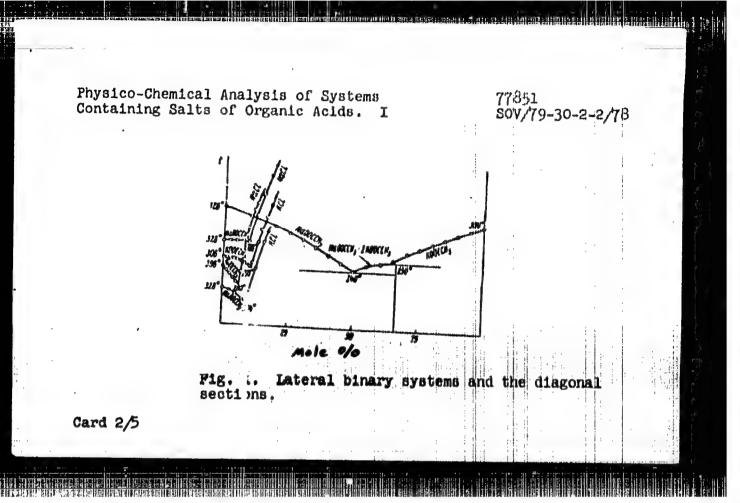
The phase diagrams for the binary systems in the ABSTRACT:

investigated mutually interacting system Na^+ , $K^+ \parallel Cl^-$, CH_3COO^- are shown in Fig. 1 (the

crystallization study was performed in test tubes

surrounded by a sleeve heater).

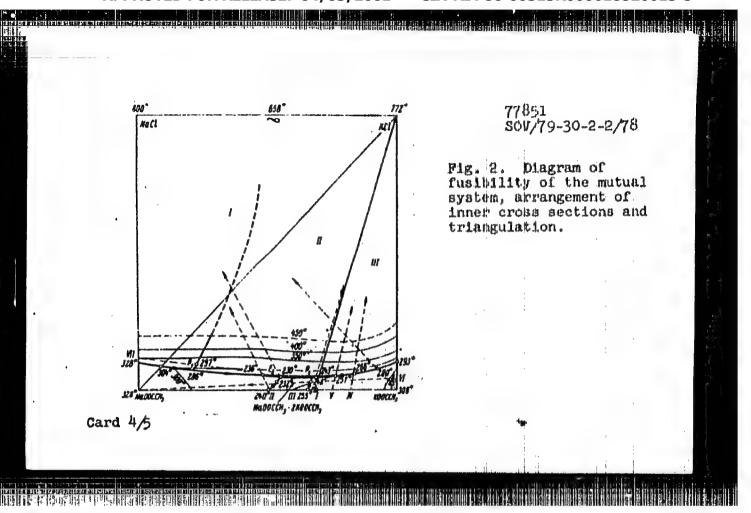
Card 1/5



Physico-Chemical Analysis of Systems Containing Salts of Organic Acids. I 77851 \$0**V/**19-30-2-2**/**78

It can be seen that sodium and potassium acetates form an incongruently melting compound, NaOOCCH₃·2KOOCCH₃ with a transition point at 255° and 65% KOOCCH₃ (eutectic is at 240° and 50% KOOCCH₃); the system NaOOCCH₃ - NaCl has a eutectic point at 328° and 10% NaCl; edtectic of the system KOOCCH₃ - KCl is at 293° and 10.5% KCl. System NaCl - KCl was studied earlier (Bergman, A. G., Nikonova, I. N., Zhur, obshchey khim., 12, 460 (1942)). The crystallization surface of the mutually interacting system is shown in Fig. 2. The system Na, K | Cl, CH₃COO is a reversibly-mutual system with a triangulating diagonal KCl - NaOOCCH₃. The crystallization fields meet in three nonvariant points (see Table 3).

Card 3/5



Physico-Chemical Analysis of Systems Containing Salts of Organic Acids. I

77851 SOV/79-30-2-2/78

Key to Table 3: (1) Point; (2) temperature; (3) composition (in mole %); (4) equilibrium phases.

	1	(3)			
(1)	(2)	Naci	Kooccij	Неооссіц	(4)
$egin{array}{c} P_1 \ E_3 \ P_0 \end{array}$	286° 230 243	7.5 5.0 5.0	22.0 55.0 86.0	40,0	NaCIKCINaOlicella KCINaOlicella KCINaOlicella NaOlicella - 2KOOCElla - KOOCElla NaOlicella - 2KOOCElla - KOOCElla

There are 2 figures; 3 tables; and 13 Soulet references.

ASSOCIATION:

Rostov-on-Don Civil-Engineering Institute (Rostovskiy-na-Donu inzhenerno-stroitel'nyy institut)
January 29, 1959

SUBMITTED:

Card 5/5

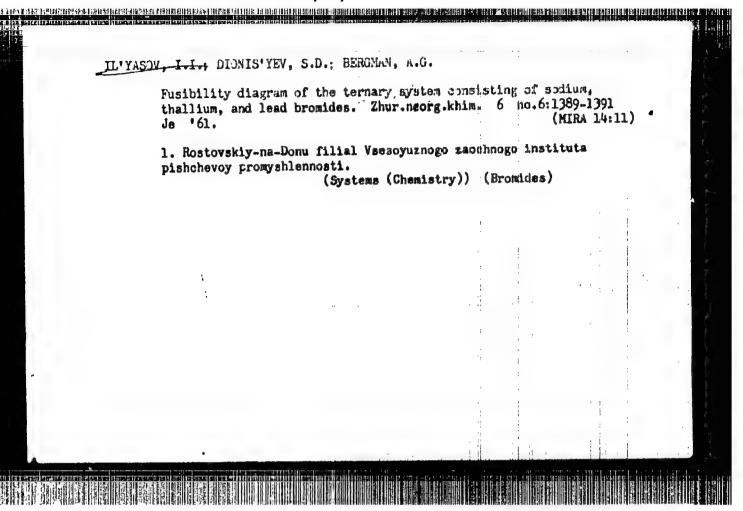
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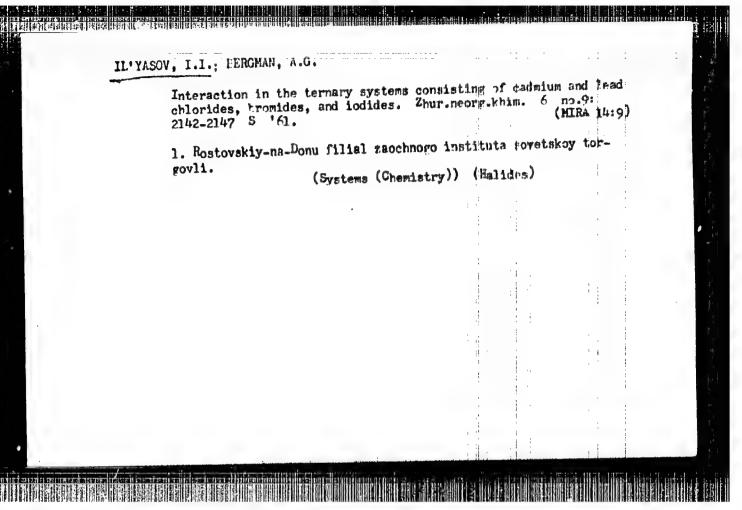
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IL!YASOV, I.I., SHCHEMELEVA, G.G., BERGMAN, A.G.

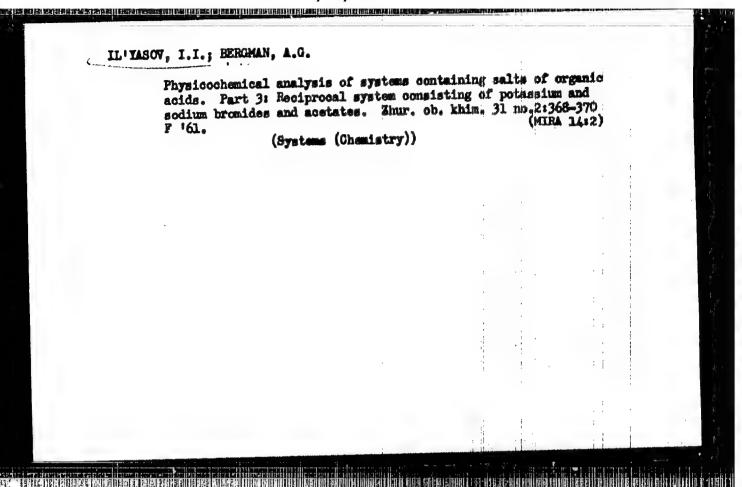
Fusibility of a ternary system of sodium, potassium, and thallium iodides. Zhur. neorg. khim. 6 no.3:699-701 Mr '61. (MIRA 14:3)

1. Rostovskiy-na-Donu filial Vsesoyuznogo zaomlnogo instituta pishehevoy promyshlennosti. (Sodium iodide) (Potassium iodide) (Thallium iodide)

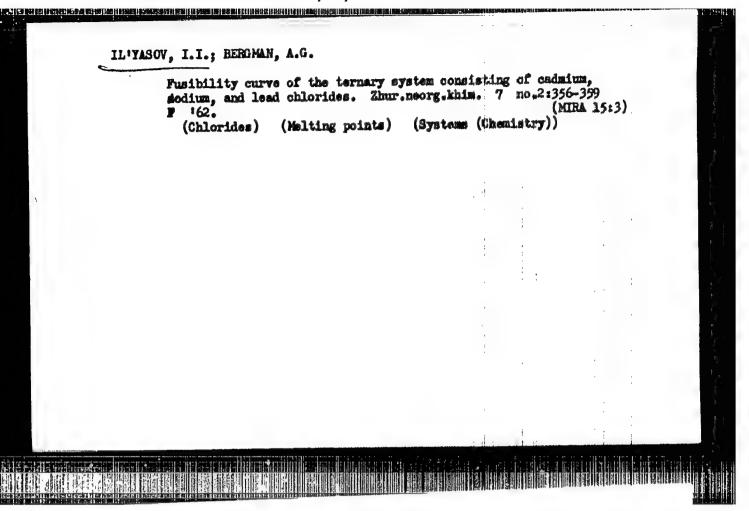


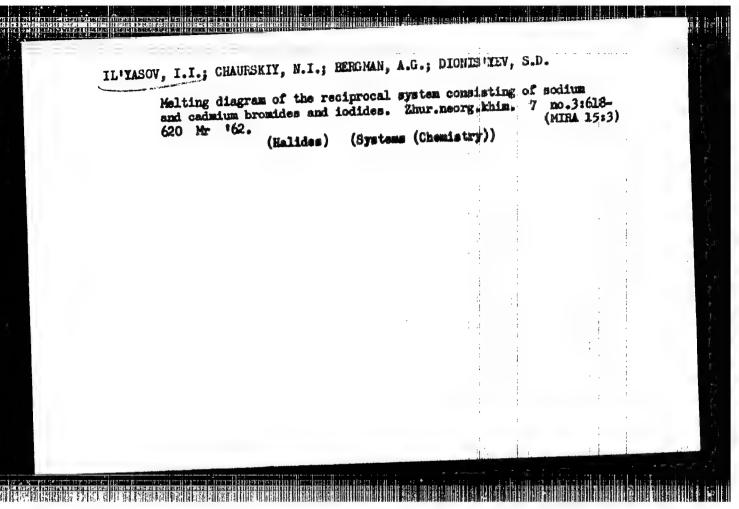


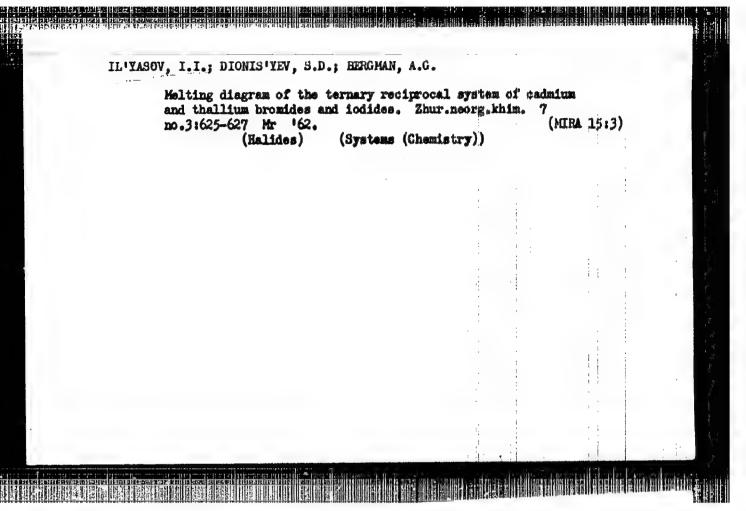
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4.1169 s/078/62/007/003/016/019 B110/B138

11.4100 AUTHORS:

Il'yasov, I. I., Bergman, A. G.

TITLE:

Pusibility of ternary systems of sodium, potassium, and

cesium chlorides

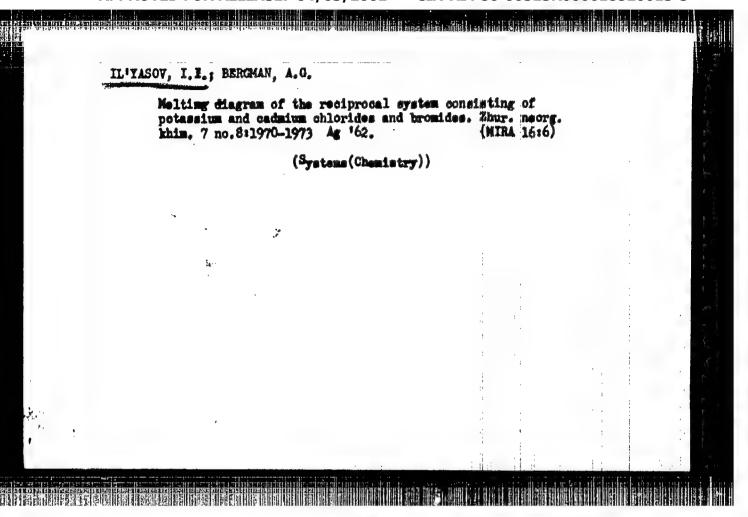
Zhurnal neorganicheskoy khimii, v. 7, no. 5, 1962, 695 - 696 PERIODICAL:

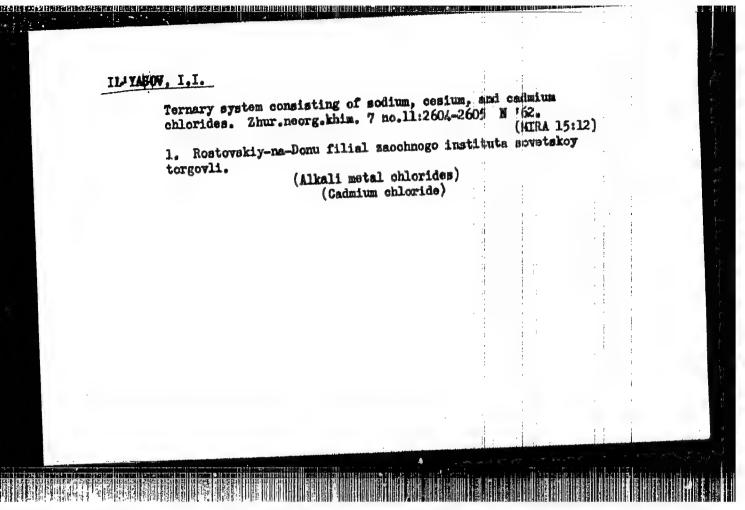
TEXT: According to its position in the periodic system the properties and salt structure of cesium differ more widely from those of sodium than potassium (to whose subgroup it belongs). The great difference in ionic

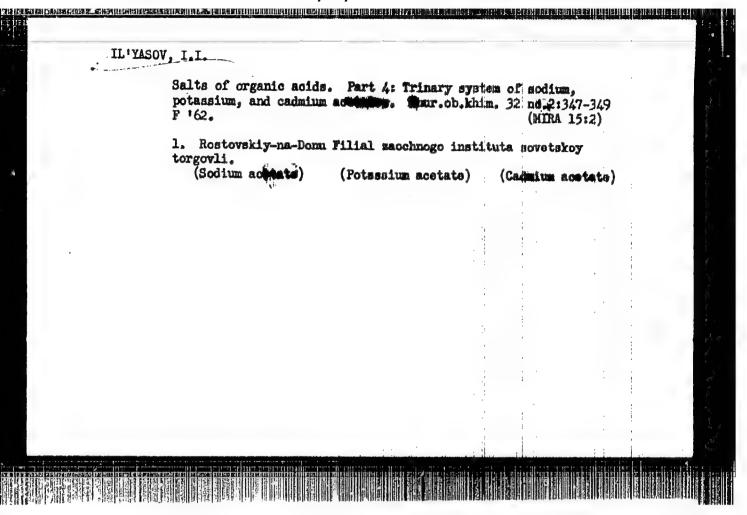
radii (K = 1.33 Å, Cs = 1.65 Å), does not, however, prevent the formation of continuous solid solutions. The authors used their own visual-polythermal method (Zh. obshch. khimii, 26, 981 (1956)). Data are expressed in moles per cent. As to the NaCl-CaCl binary systems, the authors! data on the eutectic at 34% NaCl and 493°C coincide with those of on the sursuit at 14% had and 477 (1909), but the S. F. Zhemchuzhnyy, F. Rambakh (ZhRFKhO, 41, 1785 (1909)), but the fusibility curves are somewhat more ourved. In CaCl-KCl, a continuous series of solid solutions was found with a minimum at 606°C and 36% KCl. As to NaCl-ECl, A. G. Bergman and N. M. Selivanova (Izv. Sektora fiz.-khim.

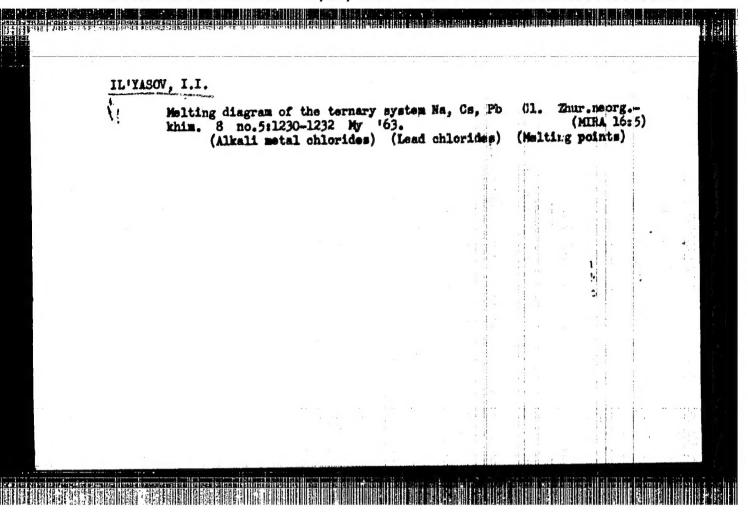
Card 1/32

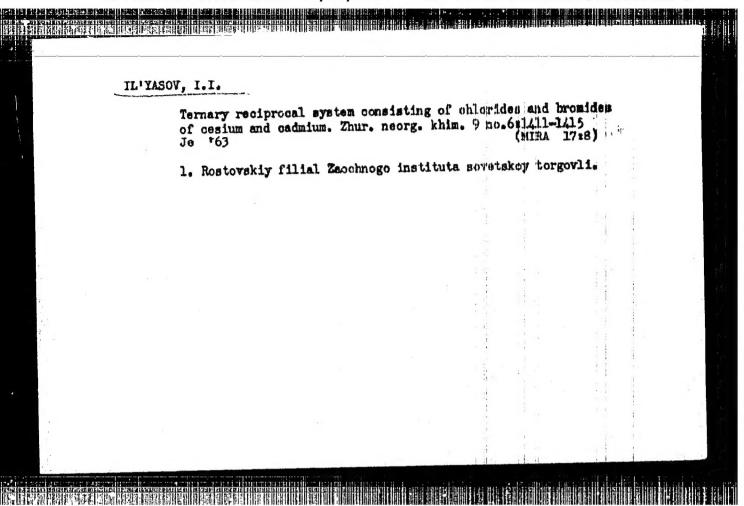
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IL'YASOV, I.I.; BERGMAN, A.G.

Ternary resilprocal systems of the halides of sodium, potassium, and cadmium. Zhur. neorg. khim. 9 no.6: 1416-1422 Je 163 (MIRA 17:8)

1. Rostovskiy filial zaochmego instituta sovetskoy torgovli.

S/079/65/035/001/002/023 D205/D507

AUTHORS:

Il'yasov, I. I., Palobekov, A. G. and Bergman, A. G.

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TITLE:

Interactions in the ternary system urea-phenol-ben-

zoic acid

PERIODICAL: Zhurnal obshchey khimii, v. 33, no. 1, 1963, 19-22

TEXT: The present work was undertaken in an effort to study systematically the interactions of urea with organic compounds. Fure materials and visual-polythermal methods were used. Melting point measurements in the binary systems phenol-urea (I), phenol-benzoic acid (II), and urea-benzoic acid (III) showed the existence of: I - a eutectic at 35°C and 6.5 mol% urea, and a transition point at 60.6°C and 33.0% CO(NH₂), corresponding to a compound CO(NH₂)₂. C₆H₅OH; II - a eutectic at 28°C and 14% C₆H₅COOH; III - a congruent melting compound 3CO(NH₂)₂. C₆H₅COOH, separating from the liquid phase at 110°C, and 2 eutectic points at 75.5°C/49.5% urea

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